



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 4, 2017 – 12:47 AM EDT

PDB ID : 2C4F
Title : crystal structure of factor VII.stf complexed with pd0297121
Authors : Kohrt, J.T.; Zhang, E.
Deposited on : unknown
Resolution : 1.72 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

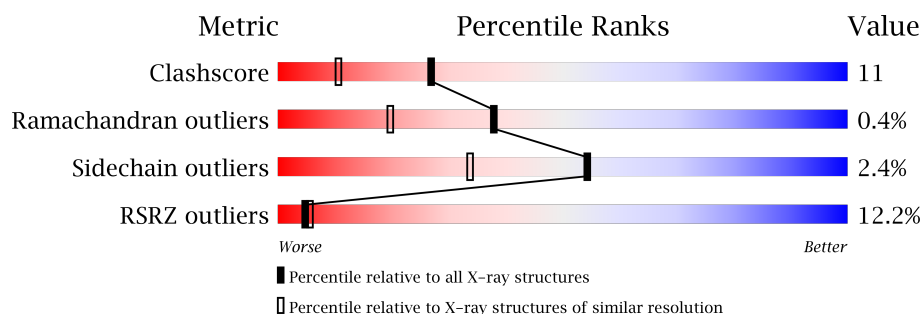
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	4841 (1.74-1.70)
Ramachandran outliers	110173	4771 (1.74-1.70)
Sidechain outliers	110143	4771 (1.74-1.70)
RSRZ outliers	101464	4426 (1.74-1.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	H	254	<div> <div>10%</div> <div>76%</div> <div>22%</div> <div>.</div> </div>
2	L	142	<div> <div>16%</div> <div>77%</div> <div>20%</div> <div>..</div> </div>
3	T	75	<div> <div>%</div> <div>87%</div> <div>13%</div> </div>
4	U	116	<div> <div>17%</div> <div>75%</div> <div>25%</div> </div>

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 5058 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called COAGULATION FACTOR VII PRECURSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	254	Total	C	N	O	S	0	0	0
			1974	1253	351	357	13			

- Molecule 2 is a protein called COAGULATION FACTOR VII PRECURSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	139	Total	C	N	O	S	0	0	0
			1117	673	185	244	15			

- Molecule 3 is a protein called TISSUE FACTOR PRECURSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	75	Total	C	N	O	S	0	0	0
			615	397	95	121	2			

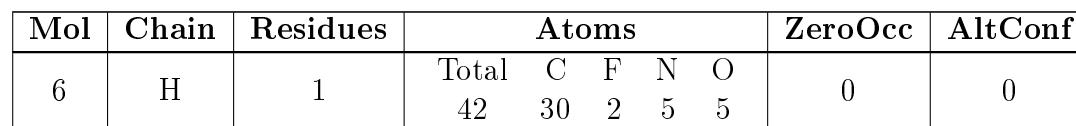
- Molecule 4 is a protein called TISSUE FACTOR PRECURSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	U	116	Total	C	N	O	S	0	0	0
			940	593	155	189	3			

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	1	Total	Ca	0	0
			1	1		
5	L	4	Total	Ca	0	0
			4	4		

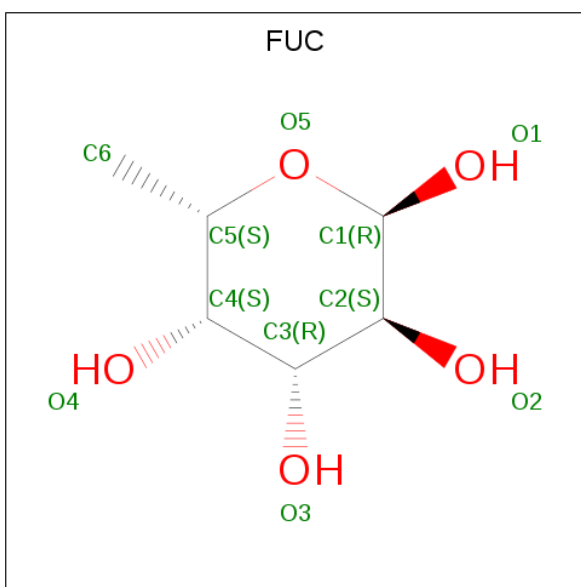
- Molecule 6 is 2-{[6-{3-[AMINO(IMINO)METHYL]PHENOXY}-4-(DIISOPROPYLAMINO)-3,5-DIFLUOROPYRIDIN-2-YL]OXY}-5-[(ISOBUTYLAMINO)CARBONYL]BENZOIC ACID (three-letter code: GIL) (formula: C₃₀H₃₅F₂N₅O₅).



- GLC
-
- Chemical structure of D-glucose (GLC) in its cyclic form (pyranose ring). The structure shows the ring carbons labeled C1(S), C2(R), C3(S), C4(S), and C5(R). The hydroxyl groups (OH) are attached to the ring carbons, and the C6 carbon is shown as a CH₂OH group. The stereochemistry is indicated by wedges and dashes: C1(S) has OH dashed; C2(R) has OH dashed; C3(S) has OH wedged; C4(S) has OH dashed; C5(R) has OH wedged. The C6 carbon is bonded to two hydroxyl groups (OH) in red, one wedged and one dashed. The C6 carbon is also bonded to an oxygen atom (O) in red, which is part of the ring. The C6 carbon is also bonded to a hydrogen atom (H) in red, which is part of the ring. The C6 carbon is also bonded to a hydrogen atom (H) in red, which is part of the ring.

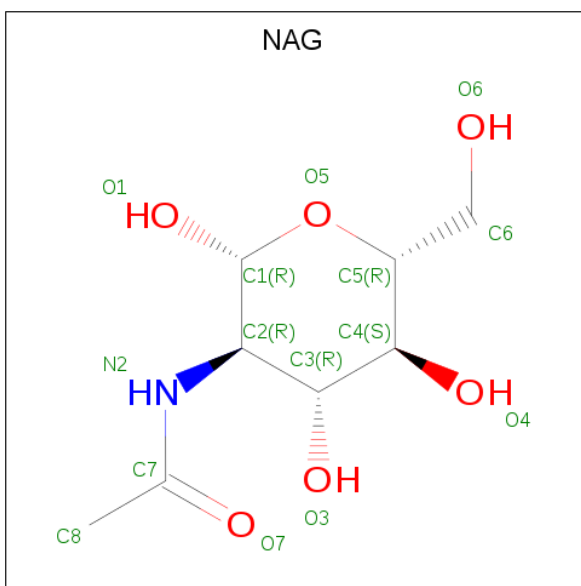
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	L	1	Total	C	O	0	0
			12	6	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	L	1	Total	C	O	0	0
			11	6	5		

- Molecule 9 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	U	1	Total	C	N	O	0	0
			14	8	1	5		

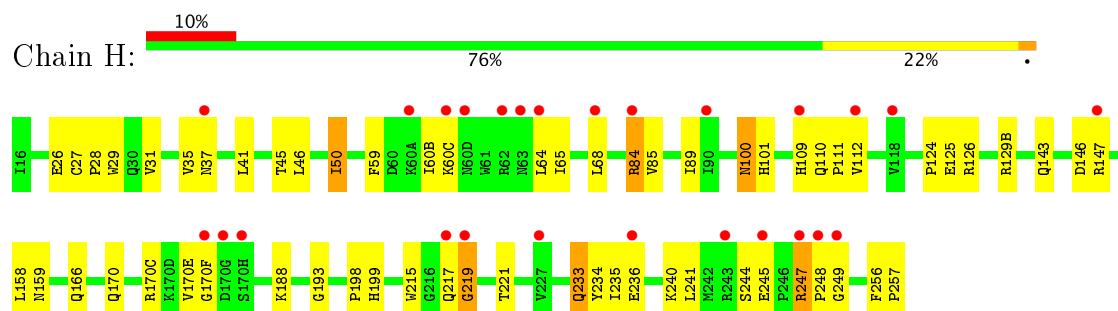
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	H	154	Total 154	O 154	0	0
10	L	63	Total 63	O 63	0	0
10	T	58	Total 58	O 58	0	0
10	U	53	Total 53	O 53	0	0

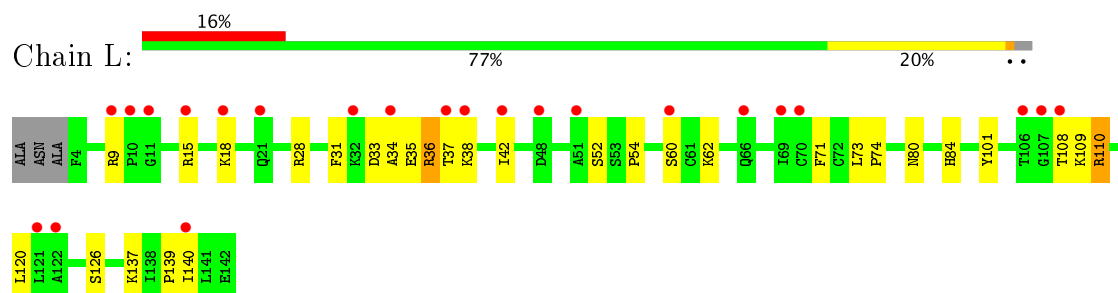
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

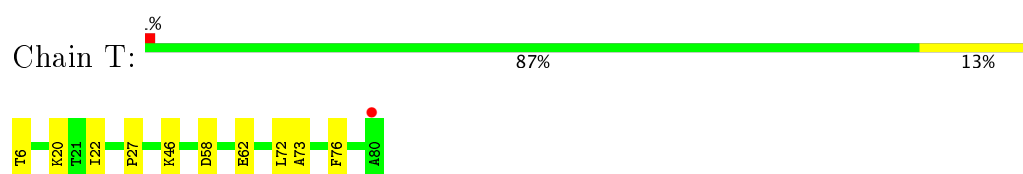
• Molecule 1: COAGULATION FACTOR VII PRECURSOR



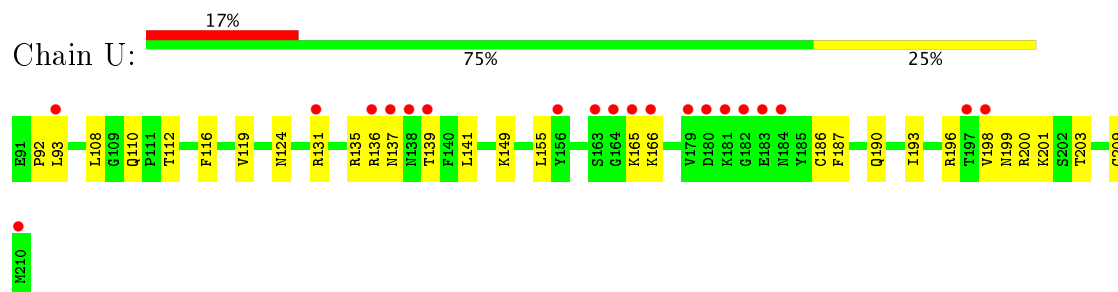
• Molecule 2: COAGULATION FACTOR VII PRECURSOR



• Molecule 3: TISSUE FACTOR PRECURSOR



• Molecule 4: TISSUE FACTOR PRECURSOR



GLOBAL-STATISTICS INFOmissingINFO

4 Model quality [i](#)

4.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, CA, GLC, FUC, GIL, CGU

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	H	0.27	0/2024	0.65	1/2755 (0.0%)
2	L	0.26	0/1011	0.53	0/1349
3	T	0.29	0/631	0.68	1/860 (0.1%)
4	U	0.27	0/958	0.63	0/1299
All	All	0.27	0/4624	0.63	2/6263 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	199	HIS	N-CA-C	-6.55	93.30	111.00
3	T	20	LYS	N-CA-C	-5.07	97.31	111.00

There are no chirality outliers.

There are no planarity outliers.

4.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	1974	0	1950	48	0
2	L	1117	0	970	34	0
3	T	615	0	598	9	0
4	U	940	0	911	25	0
5	H	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	L	4	0	0	0	0
6	H	42	0	33	3	0
7	L	12	0	12	2	0
8	L	11	0	12	4	0
9	U	14	0	13	1	0
10	H	154	0	0	3	0
10	L	63	0	0	3	0
10	T	58	0	0	2	0
10	U	53	0	0	4	0
All	All	5058	0	4499	106	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (106) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:68:LEU:HD12	1:H:112:VAL:HG11	1.37	1.04
2:L:60:SER:HB3	8:L:203:FUC:O1	1.59	1.01
2:L:74:PRO:HA	4:U:135:ARG:HH12	1.38	0.87
2:L:31:PHE:HB2	2:L:37:THR:HG22	1.61	0.81
2:L:60:SER:HB3	8:L:203:FUC:C1	2.10	0.81
2:L:60:SER:OG	2:L:71:PHE:HB2	1.83	0.78
2:L:52:SER:OG	7:L:202:GLC:H1	1.83	0.78
1:H:129(B):ARG:HD3	10:L:305:HOH:O	1.82	0.77
1:H:126:ARG:HA	1:H:235:ILE:HD12	1.66	0.77
1:H:50:ILE:HG22	1:H:111:PRO:HB3	1.68	0.76
4:U:136:ARG:O	4:U:139:THR:HG22	1.87	0.74
3:T:6:THR:HG22	4:U:93:LEU:HD22	1.71	0.73
1:H:143:GLN:NE2	1:H:146:ASP:O	2.22	0.72
2:L:71:PHE:CZ	4:U:131:ARG:HG3	2.27	0.70
2:L:52:SER:OG	7:L:202:GLC:C1	2.40	0.69
4:U:112:THR:HG23	10:U:418:HOH:O	1.92	0.69
2:L:38:LYS:O	2:L:42:ILE:HG13	1.93	0.68
2:L:62:LYS:HE3	10:L:308:HOH:O	1.92	0.68
2:L:74:PRO:HA	4:U:135:ARG:NH1	2.10	0.67
1:H:35:VAL:HG13	1:H:41:LEU:HD22	1.75	0.66
2:L:33:ASP:O	2:L:37:THR:HG23	1.96	0.66
1:H:126:ARG:CA	1:H:235:ILE:HD12	2.26	0.66
1:H:217:GLN:O	1:H:219:GLY:O	2.12	0.66
1:H:219:GLY:O	1:H:221:THR:HG23	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:U:149:LYS:HD2	10:U:450:HOH:O	1.96	0.64
1:H:125:GLU:HG3	2:L:101:TYR:CZ	2.34	0.63
1:H:124:PRO:O	1:H:235:ILE:HD13	1.98	0.62
1:H:35:VAL:CG1	1:H:41:LEU:HD22	2.29	0.62
1:H:46:LEU:HD13	1:H:68:LEU:HD11	1.82	0.61
3:T:46:LYS:HD2	10:T:124:HOH:O	2.02	0.60
1:H:147:ARG:HG2	1:H:147:ARG:O	2.01	0.60
2:L:18:LYS:HE3	10:U:408:HOH:O	2.03	0.58
1:H:64:LEU:HB3	1:H:85:VAL:HB	1.85	0.58
2:L:31:PHE:HB2	2:L:37:THR:CG2	2.31	0.58
2:L:34:ALA:O	2:L:38:LYS:HG2	2.04	0.58
1:H:26:GLU:HG3	2:L:140:ILE:HD11	1.87	0.57
1:H:166:GLN:O	1:H:170:GLN:HG3	2.05	0.57
1:H:45:THR:OG1	1:H:198:PRO:HB3	2.04	0.57
2:L:28:ARG:HA	2:L:37:THR:HG21	1.88	0.56
4:U:119:VAL:HG23	4:U:119:VAL:O	2.05	0.56
3:T:46:LYS:HD3	3:T:62:GLU:CD	2.28	0.55
1:H:170(E):VAL:HG21	1:H:217:GLN:OE1	2.07	0.54
1:H:240:LYS:HD2	10:H:504:HOH:O	2.08	0.53
4:U:198:VAL:HG22	10:U:404:HOH:O	2.09	0.52
1:H:64:LEU:HD23	1:H:85:VAL:HG11	1.91	0.52
1:H:126:ARG:NH2	1:H:233:GLN:NE2	2.58	0.51
1:H:60(C):LYS:HG2	10:H:534:HOH:O	2.10	0.51
2:L:28:ARG:HA	2:L:37:THR:CG2	2.40	0.51
1:H:84:ARG:N	1:H:84:ARG:HD2	2.25	0.51
3:T:76:PHE:CD1	4:U:92:PRO:HG2	2.46	0.50
3:T:72:LEU:HD13	3:T:73:ALA:N	2.27	0.50
1:H:126:ARG:HH22	1:H:233:GLN:NE2	2.11	0.49
4:U:166:LYS:HD2	4:U:166:LYS:N	2.28	0.49
1:H:193:GLY:HA3	6:H:302:GIL:H282	1.95	0.48
4:U:155:LEU:HD11	4:U:187:PHE:HB3	1.95	0.48
1:H:236:GLU:HG3	10:H:489:HOH:O	2.14	0.48
4:U:165:LYS:C	4:U:166:LYS:HD2	2.35	0.48
1:H:215:TRP:HB3	6:H:302:GIL:C18	2.44	0.47
1:H:215:TRP:HB3	6:H:302:GIL:H182	1.96	0.47
2:L:137:LYS:O	2:L:139:PRO:HD3	2.15	0.47
2:L:73:LEU:HD21	8:L:203:FUC:H4	1.97	0.46
3:T:22:ILE:HD13	3:T:58:ASP:HA	1.96	0.46
4:U:136:ARG:HE	4:U:141:LEU:HD21	1.79	0.46
4:U:196:ARG:O	4:U:200:ARG:NH1	2.48	0.46
1:H:158:LEU:HD11	1:H:188:LYS:HD3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:170(C):ARG:HG2	1:H:170(C):ARG:HH11	1.79	0.46
4:U:190:GLN:OE1	4:U:201:LYS:HE2	2.16	0.46
1:H:89:ILE:HG21	1:H:241:LEU:HD13	1.98	0.46
2:L:33:ASP:CG	2:L:36:ARG:HD3	2.36	0.46
4:U:136:ARG:HG2	4:U:136:ARG:HH11	1.81	0.46
1:H:35:VAL:O	1:H:35:VAL:HG23	2.16	0.46
1:H:219:GLY:C	1:H:221:THR:HG23	2.37	0.45
3:T:76:PHE:HD1	4:U:92:PRO:HG2	1.80	0.45
3:T:46:LYS:HD3	3:T:62:GLU:OE2	2.16	0.45
4:U:136:ARG:O	4:U:137:ASN:HB2	2.15	0.45
2:L:60:SER:CB	8:L:203:FUC:O1	2.47	0.45
1:H:256:PHE:HA	1:H:257:PRO:C	2.37	0.45
2:L:71:PHE:CE2	4:U:131:ARG:HG3	2.51	0.45
1:H:35:VAL:HG12	1:H:64:LEU:CD1	2.47	0.43
4:U:110:GLN:HB2	4:U:203:THR:CG2	2.48	0.43
2:L:84:HIS:HE1	10:L:345:HOH:O	2.00	0.43
1:H:125:GLU:HG3	2:L:101:TYR:CE2	2.53	0.43
2:L:110:ARG:C	2:L:110:ARG:HD2	2.39	0.43
4:U:186:CYS:HA	4:U:209:CYS:HA	2.01	0.43
3:T:46:LYS:NZ	10:T:104:HOH:O	2.53	0.42
1:H:247:ARG:HB3	1:H:248:PRO:HD2	2.01	0.42
2:L:108:THR:OG1	2:L:109:LYS:N	2.52	0.42
1:H:241:LEU:HA	1:H:244:SER:OG	2.19	0.42
1:H:27:CYS:N	1:H:28:PRO:CD	2.82	0.42
2:L:33:ASP:HB3	2:L:36:ARG:CG	2.50	0.42
4:U:124:ASN:CG	9:U:301:NAG:H82	2.39	0.42
2:L:9:ARG:CZ	2:L:15:ARG:HD3	2.50	0.42
1:H:65:ILE:HD11	1:H:84:ARG:HH12	1.84	0.42
2:L:120:LEU:HD12	2:L:126:SER:O	2.20	0.42
4:U:108:LEU:HD11	4:U:193:ILE:HG13	2.01	0.41
1:H:101:HIS:HA	1:H:234:TYR:OH	2.20	0.41
1:H:59:PHE:HA	1:H:60(B):ILE:HG12	2.02	0.41
2:L:108:THR:HG23	2:L:109:LYS:N	2.35	0.41
1:H:245:GLU:O	1:H:247:ARG:NE	2.53	0.41
2:L:35:CGU:O	2:L:38:LYS:HG2	2.20	0.41
1:H:248:PRO:HG2	1:H:249:GLY:H	1.86	0.41
1:H:31:VAL:HG22	1:H:68:LEU:HD21	2.03	0.41
4:U:116:PHE:HA	4:U:124:ASN:O	2.21	0.41
2:L:54:PRO:O	2:L:80:ASN:HB3	2.21	0.40
1:H:100:ASN:HD22	1:H:100:ASN:HA	1.61	0.40
1:H:109:HIS:ND1	1:H:110:GLN:HG2	2.37	0.40

There are no symmetry-related clashes.

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	252/254 (99%)	238 (94%)	12 (5%)	2 (1%)	22	7
2	L	127/142 (89%)	122 (96%)	5 (4%)	0	100	100
3	T	73/75 (97%)	72 (99%)	1 (1%)	0	100	100
4	U	112/116 (97%)	107 (96%)	5 (4%)	0	100	100
All	All	564/587 (96%)	539 (96%)	23 (4%)	2 (0%)	38	20

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	219	GLY
1	H	170(F)	GLY

4.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	216/216 (100%)	208 (96%)	8 (4%)	39	17
2	L	113/114 (99%)	111 (98%)	2 (2%)	64	46
3	T	70/70 (100%)	69 (99%)	1 (1%)	71	57
4	U	109/109 (100%)	108 (99%)	1 (1%)	82	74
All	All	508/509 (100%)	496 (98%)	12 (2%)	54	34

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	29	TRP
1	H	37	ASN
1	H	50	ILE
1	H	84	ARG
1	H	100	ASN
1	H	159	ASN
1	H	233	GLN
1	H	247	ARG
2	L	36	ARG
2	L	110	ARG
3	T	27	PRO
4	U	199	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	37	ASN
1	H	63	ASN
1	H	100	ASN
1	H	110	GLN
1	H	143	GLN
1	H	159	ASN
1	H	202	HIS
1	H	233	GLN
1	H	239	GLN
2	L	80	ASN
2	L	84	HIS
3	T	31	ASN
4	U	171	ASN
4	U	199	ASN

4.3.3 RNA ⓘ

There are no RNA molecules in this entry.

4.4 Non-standard residues in protein, DNA, RNA chains ⓘ

10 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul

statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	CGU	L	14	2,5	5,11,12	1.10	1 (20%)	4,14,16	2.06	1 (25%)
2	CGU	L	16	2,5	5,11,12	1.19	1 (20%)	4,14,16	2.11	1 (25%)
2	CGU	L	19	2,5	5,11,12	1.13	1 (20%)	4,14,16	1.98	1 (25%)
2	CGU	L	20	2	5,11,12	1.38	1 (20%)	4,14,16	1.99	1 (25%)
2	CGU	L	25	2,5	5,11,12	0.98	0	4,14,16	2.04	1 (25%)
2	CGU	L	26	2,5	5,11,12	1.05	1 (20%)	4,14,16	2.15	1 (25%)
2	CGU	L	29	2,5	5,11,12	1.01	1 (20%)	4,14,16	1.90	1 (25%)
2	CGU	L	35	2	5,11,12	1.02	1 (20%)	4,14,16	2.08	1 (25%)
2	CGU	L	6	2	5,11,12	1.01	1 (20%)	4,14,16	1.91	1 (25%)
2	CGU	L	7	2	5,11,12	1.22	1 (20%)	4,14,16	2.11	1 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CGU	L	14	2,5	-	0/4/14/16	0/0/0/0
2	CGU	L	16	2,5	-	0/4/14/16	0/0/0/0
2	CGU	L	19	2,5	-	0/4/14/16	0/0/0/0
2	CGU	L	20	2	-	0/4/14/16	0/0/0/0
2	CGU	L	25	2,5	-	0/4/14/16	0/0/0/0
2	CGU	L	26	2,5	-	0/4/14/16	0/0/0/0
2	CGU	L	29	2,5	-	0/4/14/16	0/0/0/0
2	CGU	L	35	2	-	0/4/14/16	0/0/0/0
2	CGU	L	6	2	-	0/4/14/16	0/0/0/0
2	CGU	L	7	2	-	0/4/14/16	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	6	CGU	CA-C	2.00	1.52	1.50
2	L	29	CGU	CA-C	2.04	1.52	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	35	CGU	CA-C	2.07	1.53	1.50
2	L	26	CGU	CA-C	2.16	1.53	1.50
2	L	14	CGU	CA-C	2.25	1.53	1.50
2	L	19	CGU	CA-C	2.31	1.53	1.50
2	L	16	CGU	CA-C	2.50	1.53	1.50
2	L	7	CGU	CA-C	2.54	1.53	1.50
2	L	20	CGU	CA-C	2.93	1.54	1.50

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	26	CGU	CB-CG-CD2	-4.03	105.07	112.81
2	L	16	CGU	CB-CG-CD2	-3.67	105.76	112.81
2	L	7	CGU	CB-CG-CD2	-3.51	106.06	112.81
2	L	35	CGU	CB-CG-CD2	-3.38	106.32	112.81
2	L	14	CGU	CB-CG-CD2	-3.34	106.39	112.81
2	L	29	CGU	CB-CG-CD2	-3.32	106.42	112.81
2	L	25	CGU	CB-CG-CD2	-3.29	106.49	112.81
2	L	19	CGU	CB-CG-CD2	-3.26	106.54	112.81
2	L	20	CGU	CB-CG-CD2	-3.25	106.56	112.81
2	L	6	CGU	CB-CG-CD2	-3.09	106.88	112.81

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	35	CGU	1	0

4.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

4.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 5 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GIL	H	302	-	40,44,44	1.66	9 (22%)	51,63,63	1.30	4 (7%)
7	GLC	L	202	-	12,12,12	0.26	0	17,17,17	0.33	0
8	FUC	L	203	-	11,11,11	0.31	0	15,16,16	0.41	0
9	NAG	U	301	4	14,14,15	0.53	0	15,19,21	0.69	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GIL	H	302	-	-	0/33/37/37	0/3/3/3
7	GLC	L	202	-	-	0/2/22/22	0/1/1/1
8	FUC	L	203	-	-	0/0/20/20	0/1/1/1
9	NAG	U	301	4	-	0/6/23/26	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	H	302	GIL	C6-C1	2.12	1.42	1.39
6	H	302	GIL	C22-C21	2.21	1.45	1.40
6	H	302	GIL	C2-C1	2.27	1.43	1.39
6	H	302	GIL	C13-C14	2.34	1.43	1.39
6	H	302	GIL	C11-N3	2.52	1.36	1.32
6	H	302	GIL	C23-C24	2.98	1.43	1.39
6	H	302	GIL	C13-C12	2.98	1.44	1.39
6	H	302	GIL	C15-N3	3.09	1.36	1.32
6	H	302	GIL	C12-C11	4.32	1.43	1.38

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	H	302	GIL	O1-C11-N3	-2.22	115.42	119.48
6	H	302	GIL	O2-C21-C26	-2.01	115.57	120.82
6	H	302	GIL	C17-C10-N4	2.22	117.10	112.37
6	H	302	GIL	C16-N4-C10	6.55	127.64	116.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	H	302	GIL	3	0
7	L	202	GLC	2	0
8	L	203	FUC	4	0
9	U	301	NAG	1	0

4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
4	U	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	U	158:TRP	C	163:SER	N	9.35

5 Fit of model and data ⓘ

5.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	H	254/254 (100%)	0.64	26 (10%) 7 9	10, 19, 39, 61	1 (0%)
2	L	129/142 (90%)	1.04	23 (17%) 2 2	16, 27, 44, 51	0
3	T	75/75 (100%)	0.37	1 (1%) 77 82	12, 18, 25, 30	0
4	U	116/116 (100%)	1.16	20 (17%) 2 2	15, 23, 50, 55	1 (0%)
All	All	574/587 (97%)	0.80	70 (12%) 5 5	10, 22, 41, 61	2 (0%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	U	137	ASN	8.5
4	U	181	LYS	7.8
1	H	248	PRO	7.6
4	U	182	GLY	7.2
2	L	9	ARG	6.6
4	U	138	ASN	6.5
1	H	170(G)	ASP	6.3
4	U	156	TYR	5.7
1	H	247	ARG	5.2
2	L	122	ALA	5.2
1	H	62	ARG	5.0
1	H	147	ARG	5.0
4	U	197	THR	5.0
2	L	106	THR	4.9
1	H	219	GLY	4.6
4	U	164	GLY	4.5
1	H	37	ASN	4.4
3	T	80	ALA	4.4
2	L	107	GLY	4.3
4	U	163	SER	4.3
4	U	136	ARG	4.2

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Mol	Chain	Res	Type	RSRZ
2	L	15	ARG	4.1
4	U	139	THR	3.8
1	H	249	GLY	3.8
4	U	180	ASP	3.7
2	L	21	GLN	3.6
2	L	108	THR	3.5
4	U	198	VAL	3.5
4	U	165	LYS	3.4
2	L	42	ILE	3.4
1	H	109	HIS	3.1
1	H	84	ARG	3.1
4	U	183	GLU	3.0
1	H	64	LEU	3.0
4	U	184	ASN	2.9
2	L	66	GLN	2.9
1	H	170(F)	GLY	2.9
2	L	10	PRO	2.9
2	L	18	LYS	2.9
1	H	245	GLU	2.9
1	H	60(C)	LYS	2.8
2	L	32	LYS	2.8
2	L	69	ILE	2.7
2	L	140	ILE	2.7
4	U	131	ARG	2.7
1	H	60(D)	ASN	2.7
2	L	11	GLY	2.6
1	H	68	LEU	2.6
2	L	51	ALA	2.5
1	H	243	ARG	2.5
1	H	90	ILE	2.5
4	U	166	LYS	2.5
1	H	236	GLU	2.5
2	L	37	THR	2.4
4	U	179	VAL	2.3
1	H	217	GLN	2.3
4	U	93	LEU	2.3
2	L	70	CYS	2.3
2	L	34	ALA	2.3
1	H	118	VAL	2.3
2	L	121	LEU	2.2
2	L	60	SER	2.2
4	U	210	MET	2.2

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Mol	Chain	Res	Type	RSRZ
2	L	38	LYS	2.2
1	H	63	ASN	2.2
2	L	48	ASP	2.1
1	H	112	VAL	2.1
1	H	227	VAL	2.1
1	H	60(A)	LYS	2.1
1	H	170(H)	SER	2.1

5.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CGU	L	7	12/13	0.81	0.20	-	30,33,36,38	0
2	CGU	L	19	12/13	0.75	0.33	-	46,51,54,54	0
2	CGU	L	6	12/13	0.81	0.23	-	26,34,37,38	0
2	CGU	L	14	12/13	0.71	0.29	-	37,46,48,49	0
2	CGU	L	25	12/13	0.87	0.20	-	25,26,28,29	0
2	CGU	L	29	12/13	0.90	0.13	-	25,27,30,30	0
2	CGU	L	26	12/13	0.85	0.20	-	23,29,33,35	0
2	CGU	L	16	12/13	0.86	0.17	-	34,35,37,37	0
2	CGU	L	35	12/13	0.42	0.47	-	39,51,53,54	0
2	CGU	L	20	12/13	0.72	0.34	-	44,50,52,53	0

5.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
9	NAG	U	301	14/15	0.81	0.21	1.69	24,31,41,41	0
8	FUC	L	203	11/11	0.67	0.19	0.87	35,36,39,42	0
6	GIL	H	302	42/42	0.88	0.14	0.46	14,25,34,36	0
5	CA	H	301	1/1	0.98	0.13	0.16	29,29,29,29	0
5	CA	L	201	1/1	0.97	0.05	-1.69	28,28,28,28	0
7	GLC	L	202	12/12	0.70	0.19	-	50,52,52,52	0
5	CA	L	204	1/1	0.96	0.16	-	46,46,46,46	0
5	CA	L	205	1/1	0.86	0.14	-	85,85,85,85	0
5	CA	L	206	1/1	0.95	0.14	-	45,45,45,45	0

5.5 Other polymers [i](#)

There are no such residues in this entry.